

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072123\
 Data File : BG058383.D
 Acq On : 21 Jul 2023 11:48
 Operator : CG/JU
 Sample : PB154123BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Jul 21 22:16:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 18 02:19:39 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.894	152	46581	20.000	ng/ul	0.00
20) Naphthalene-d8	10.697	136	197919	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.533	164	139052	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.277	188	348772	20.000	ng/ul	0.00
79) Chrysene-d12	21.525	240	326504	20.000	ng/ul	0.00
88) Perylene-d12	24.657	264	381139	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.335	96	8938	7.062	ng/uL	0.00
4) Pyridine-d5	3.746	84	124437	33.126	ng/ul	0.00
7) Phenol-d5	7.060	99	143146	32.395	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.218	67	92187	33.378	ng/ul	0.00
11) 2-Chlorophenol-d4	7.424	132	103372	33.420	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	103042	30.571	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	53155	35.830	ng/ul	0.00
24) 2-Nitrophenol-d4	9.774	143	59737	36.884	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.321	165	100830	32.476	ng/ul	0.00
31) 4-Chloroaniline-d4	10.832	131	145812	32.397	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	369460	34.456	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	393957	35.059	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	51229	31.635	ng/ul	0.00
60) Fluorene-d10	15.520	176	312024	35.148	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.644	200	58266	32.790	ng/ul	0.00
73) Anthracene-d10	17.377	188	540211	35.847	ng/ul	0.00
81) Pyrene-d10	19.663	212	667532	34.615	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	662273	35.951	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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